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Technical Report No. AF-105

WAVE FUNCTIONS, ENERGIES, AND REDUCED MATRIX
ELEMENTS OF THE 5f<sup>3</sup> CONFIGURATION
IN INTERMEDIATE COUPLING

Ьу

Hannah M. Crosswhite



THE JOHNS HOPKINS UNIVERSITY CARLYLE BARTON LABORATORY BALTIMORE, MD.

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ERRATA - AF 104 - A Forming Study of Point-Contact Tunnel
Diodes - H. J. Lory - May 1963

Page 11 line 9 reads are snic - should be arsenic Page 20 line 12 reads normal of N- should be - normal or N- Page 35 equation (8) should be  $\sqrt{2}$  2hs e in the denominator Page 88 line 14 reads H on D<sub>o</sub>, should read H or D<sub>o</sub>.

### THE JOHNS HOPKINS UNIVERSITY CARLYLE BARTON LABORATORY BALTIMORE, MARYLAND

Technical Report No. AF-105

### WAVE FUNCTIONS, ENERGIES, AND REDUCED MATRIX ELEMENTS OF THE 5f<sup>3</sup> CONFIGURATION IN INTERMEDIATE COUPLING

by

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### **ABSTRACT**

Since the crystal field can be considered as a perturbation of the free ion spectra of actinide ions in certain crystals, we have calculated energy levels and wave functions of the  $5f^3$  configuration, which is the ground configuration of  $U^{3+}$ , in intermediate coupling for values of  $\chi = \zeta / F_2$  of 7 to 11. Energy levels are given in units of  $F_2$ , and wave functions are expressed as linear combinations of Russell-Saunders functions. Hydrogenic and non-hydrogenic approximations are compared.

Reduced matrix elements in intermediate coupling are given so that crystal field splittings may be calculated. The splittings of the two lowest levels of  ${\tt U}^{3+}$  in a weak field with and without J-mixing are compared.

### I. INTRODUCTION

In the triply ionized actinides the optically active electrons belong to the 5f<sup>n</sup> configuration<sup>1, 2</sup>, so comparison of their spectra with those of the triply ionized rare earths, whose optically active electrons belong to the 4f<sup>n</sup> configurations, is of considerable interest. The effect of the crystal field on rare earth ions is small enough to be treated as a perturbation of the free ion spectra; this is also true for the actinide ions, at least in some crystals 3-9. In a free ion, electric dipole transitions within a configuration are forbidden, but the odd parity part of the crystal potential will mix configurations of opposite parity so that "forced" electric dipole transitions can take place as well as the allowed magnetic dipole transitions. The strength of the forced electric dipole transition depends inversely upon the energy differences between mixing configurations 10, 11. crystals containing triply ionized actinides total absorption begins at about 4000A, and it is assumed that this is due to the allowed 5f-6d transitions. (Knowledge of the free ion spectra would be useful here). Since the analogous 4f-5d transitions in rare earth crystals take place at about 2000 A<sup>12</sup>, more intense spectra can be expected from this effect in actinide crystals than in rare earth crystals.

The ground configuration of  $U^{3+}$  is  $5f^3$ . Since nothing is known experimentally about the free ion spectrum of  $U^{3+}$ , we have made calculations on the  $5f^3$  configuration to aid in the interpretation of crystal spectra.

### II. FREE ION

The Hamiltonian for the free ion is

$$H_f = H_o + H_e + H_s$$

where  $H_0$  is the interaction of the 5f electrons with the core (constant for the configuration)

 $\ensuremath{\mathrm{H}_{\mathrm{e}}}$  is the electrostatic interaction between the three 5f electrons

H<sub>s</sub> is the spin orbit interaction.

Higher order effects, including spin-spin and spin-other-orbit interactions will be neglected.

The electrostatic interaction is

$$H_e = \sum_{i>j} \frac{e^2}{r_{ij}}$$
 i, j = 1, 2, 3

where electrons in closed shells are assumed to affect the energies of the 5f electrons only by a contribution to the central potential. We can expand  $\frac{1}{r_{ij}}$  in terms of the Legendre polynomials  $^{13}$ 

$$\frac{1}{r_{ij}} = \sum_{k=0}^{\infty} \frac{r_n^n}{r_b^n} P_n(\cos \omega)$$

where  $r_a$  is the smaller and  $r_b$  the larger of the distances of  $r_i$  and  $r_j$  from the nucleus, and w is the angle between  $r_i$  and  $r_j$ . Then we can separate the radial and angular components and write

$$H_e = \sum_{n} f^n F_n$$
  $n = 2, 4, 6$  for f electrons

Matrix elements of  $f^n$  are essentially integrals of spherical harmonics over the angular wave function, and have been evaluated by Carlson  $f^n$  for  $f^n$  using Racah's method. The  $F_n$  are integrals of  $\frac{r^n}{n+1}$  over the 5f radial wave functions (Slater integrals). If 5f hydrogenic wave functions are assumed, that is, in a Coulomb field, Judd  $f^n$  has calculated

$$F_4/F_2 = 0.1422$$
 and  $F_6/F_2 = 0.0161$ 

The spin orbit interaction is

$$H_{s} = \sum_{i=1}^{3} \xi(r_{i}) \vec{l}_{i} \cdot \vec{s}_{i}$$

ζ, the integral of ξ(r<sub>i</sub>) over the radial wave function, is determined experimentally due to lack of knowledge of the wave function. The matrix elements of the angular part of H<sub>s</sub> have been calculated for f<sup>3</sup> by Judd and Loudon<sup>16</sup>. In U<sup>3+</sup> the spin orbit interaction is rather large, therefore we had to perform intermediate coupling calculations. Since J is a good quantum number, this involved diagonalizing matrices for each J value, 1/2 to 17/2. The electrostatic matrices, assuming 5f hydrogenic wave functions, and spin orbit matrices have been diagonalized simultaneously, by using a program developed

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by Wybourne\*, for various values of  $\chi = \zeta/F_2$  expressing energies in terms of  $F_2$  and wave functions as linear combinations of Russell-Saunders states. Figure 1 shows the variation in energy levels of  $f^3$  with  $\chi$ , for  $\chi=7$  to 11. The levels are named by the Russell-Saunders states into which they go at  $\zeta=0$ . For  $F_2=200$ ,  $\chi=8$ , (the order of magnitude appropriate for  $U^{3+})^{17}$ , it is noted that the configuration extends 35,000 cm<sup>-1</sup>. In comparison, the Nd<sup>3+</sup> 4f<sup>3</sup> configuration extends 67,000 cm<sup>-1</sup>. The 41 levels are each (2J+1)-fold degenerate.

The use of the 5f hydrogenic wave functions for U<sup>3+</sup> is questionable. Lammermann and Conway<sup>8</sup> feel that the Coulomb approximation is satisfactory for the interpretation of the Pu<sup>3+</sup>(5f<sup>5</sup>) spectrum. Cohen has done a relativistic self-consistent calculation for the normal uranium ion, and Winocur has calculated the following set of Slater integrals from Cohen's eigenfunctions<sup>18</sup>

$$F_4/F_2 = 0.159$$
  $F_6/F_2 = 0.0204$ .

The discrepancy between a similar calculation for PrIV and the experimental values indicates that these should be treated with reserve <sup>17</sup>.

McLaughlin determined  $F_2$ ,  $F_4$ , and  $F_6$  experimentally by a least squares fit to the spectrum of  $UCl_4(5f^2)$  and found the ratios to be

<sup>\*</sup> Private communication.

$$F_4/F_2 = 0.1468$$
  $F_6/F_2 = 0.0219$ 

We have calculated energy levels and wave functions of  $f^3$  using these last ratios with  $\chi=8.0$ , and the energy levels are compared with hydrogenic ones in Figure 2.  $F_2$  has been chosen to put the  $^4I_{11/2}$  level at 4500 cm $^{-1}$  from the ground state where it has been observed  $^{19,20}$ . It is apparent that the position of the  $^4F_{3/2}$  level relative to the  $^4I_{13/2}$  is depressed in the non-hydrogenic case, and above 10,000 cm $^{-1}$ , where the energy levels become more dense, the grouping is much different in the two cases.

### III. CRYSTAL FIELD

We are interested in crystals in which the splitting of the J manifold is small compared to the spin-orbit splitting, so in a first approximation the crystal field interaction can be treated as a perturbation of the free ion energy levels. However, since the 5f wave functions are more extensive than the 4f wave functions <sup>22</sup>, somewhat larger crystal field effects can be expected in the actinides than in the rare earths. Each J level of an ion with an odd number of electrons will split in a crystal field into (2J + 1)/2 components at the most due to Kramers degeneracy.

 $\begin{tabular}{ll} The Hamiltonian for the crystal field interaction, assuming \\ a purely electrostatic field is $^{23}$ \\ \end{tabular}$ 

$$H_c = \sum_{kq}^{q} A_k^q \langle r^k \rangle V_k^q(\theta, \phi) |q| \le k, k = 2, 4, 6 \text{ for f electrons.}$$

The number of terms in the expansion depends on the symmetry of the ion site in the crystal. The coefficients  $A_k^q$  depend on the lattice, and, although it is theoretically possible to calculate them if the positions and charge distributions of the ions in the crystal are known, results have not been satisfactory. Therefore,  $A_k^q \langle r^k \rangle$  are customarily treated as experimentally determined parameters. The  $V_k^q$  are normalized associate spherical harmonics which transform in the same manner as the unit tensor operators  $U_k^q$  and are related to them by constant factors. We may use either set for calculations. Matrix elements of the crystal potential, for LS coupling then, can be obtained by using the relation  $^{45}$ 

 $\langle \texttt{a} \; \texttt{LSJM} \big| \texttt{U}_q^k \big| \texttt{a'L'S'J'M'} \rangle = \delta(\texttt{SS'}) (-1)^{J-M} \binom{J-kJ'}{-M \cdot q \cdot M'} \\ \langle \texttt{a} \; \texttt{LSJ} \; \big| \; \big| \texttt{U}^k \big| \; \big| \texttt{a'L'S'J'} \rangle,$ 

where the reduced matrix element is

 $\langle a LSJ \| U^k \| a'L'S'J' \rangle = (-1)^{S+k+L'+J} \Big[ (2J+1)(2J'+1) \Big]^{1/2} \left\{ \begin{matrix} L \ J \ S \\ J' \ L' \ k \end{matrix} \right\} \langle a LS \| U^k \| a'L'S \rangle.$   $3-j \ symbols \left( \begin{matrix} J \ kJ' \\ -Mq \ M' \end{matrix} \right) \ and \ 6-j \ symbols \left\{ \begin{matrix} L \ J \ S \\ J' \ L' \ k \end{matrix} \right\} \ are \ vector \ coupling$  25  $coefficients \ and \ are \ available \ in \ tabular \ form^{24}. \ Judd \ has \ tabulated$   $the \ double \ reduced \ matrix \ elements \ \langle a LS \| U^k \| a'L'S \rangle \ for \ f^3. \ If \ J$   $mixing \ is \ neglected, \ the \ matrix \ elements \ may \ be \ more \ simply$   $calculated \ from$ 

$$\langle \alpha \text{LSJM} \big| \, V_k^{\mathbf{q}} \big| \alpha' \text{L'SJM'} \rangle = \langle \alpha \text{LSJ} \big| \big| \, V_k^{\phantom{\dagger}} \big| \big| \alpha' \text{L'SJ} \rangle \,\, f_k^{\phantom{\dagger}} S_k^{\mathbf{q}}(M),$$

where  $S_k^q(M)$  are the Stevens coefficients  $^{26}$  and the  $f_k$ , which depend only on J, are proportionality constants chosen to make the  $S_k^o$  integers.

 $U^{3+}$  states can be represented by linear combinations of Russell-Saunders states with coefficients depending on  $\chi$ ; therefore the crystal field matrix elements will also depend on  $\chi$ , and we have obtained them for  $\Delta J$  = 0 by the same transformation which diagonalized the  $5f^3$  energy matrix.

Table I presents the results of calculations we have made on the  $5f^3$  configuration using hydrogenic wave functions. For each J value for  $\chi=7.0$  to 11.0, energies in units of  $F_2$  are given in the first row. In the columns below each energy level are the coefficients

COLX

of the LS wave functions for that level. Reduced matrix elements in intermediate coupling are given in terms of  $\langle aLSJ || V_k || aL'SJ \rangle$   $f_k \times 10^3$  in order that they may be used with the Stevens coefficients given in Table II<sup>26, 27</sup>. The last row of each column gives the Landé g-value of the level in intermediate coupling.

As an example of the use of the tables, the splitting of a  $^4F_{5/2}$  level in  $D_{3h}$  symmetry for  $\chi$  = 8 is (-0.00937)  $S_2^O(M)$  + (-0.02391)  $S_4^O(M)$  or

 $(-0.00937) \times 5 \times A_2^{o} + (-0.02391) \times 1 \times A_4^{o} = -0.04685 A_2^{o} - 0.02391 A_4^{o}, m = 5/2$   $-(0.00937) \times (-1) \times A_2^{o} - (0.02391) \times (-3) \times A_4^{o} = 0.00937 A_2^{o} + 0.07173 A_4^{o}, m = 3/2$   $-(0.00937) \times (-4) \times A_2^{o} - (0.02391) \times 2 A_4^{o} = 0.03748 A_2^{o} - 0.04782 A_4^{o}, m = 1/2 .$ 

In a crystal field J is only approximately a good quantum number, so that J levels will also be mixed. A complete J mixing calculation would involve diagonalizing three 60- and 61-dimensional matrices, and does not seem to be worthwhile until more is known about the wave functions. However, in order to get an idea of the magnitude involved, J mixing of the  ${}^4I_{9/2}$  and  ${}^4I_{11/2}$  states can be calculated by second order perturbation theory. (For higher levels which become more nearly degenerate this cannot be done).

We have calculated the splittings of the  $^4I_{9/2}$  and  $^4I_{11/2}$  states using 5f hydrogenic wave functions for  $\chi$  = 8.0 and the fictitious crystal field parameters

CATW

$$A_2^o \langle \mathbf{r}^2 \rangle = 170$$
  $A_6^o \langle \mathbf{r}^6 \rangle = -100$   $A_4^o \langle \mathbf{r}^4 \rangle = -80$   $A_6^o \langle \mathbf{r}^6 \rangle = 1100$ 

which are close to those found by Gruber  $^4$  for Am  $^{3+}$  in LaCl $_3$ . The effect of the  $^4I_{11/2}$  state on the  $^4I_{9/2}$  state 4500 cm  $^{-1}$  below it is to depress the ground Stark level by 27 cm  $^{-1}$  and to increase slightly the overall splitting. The calculated splittings of the  $^4I_{11/2}$  and  $^4I_{9/2}$  levels using the above parameters, with and without J mixing, are shown in Figure 3.

PATW

Experimentally, a complication occurs in uranium doped crystals in that uranium readily ionizes to U  $^{4+}$  and less easily to U  $^{3+}$  so that crystals nominally containing U  $^{3+}$  sometimes contain both ions, as was the case for CaF  $_2$ :U  $^{3+}$ .  $^{28,29}$  The ground configuration of U  $^{4+}$  is 5f  $^2$ . In Figure 4 we compare McLaughlin's energy level diagram for U  $^{4+}$  with our calculated one for 5f  $^3$  for reasonable values of F  $_2$  and  $\chi$ , and find that the energy levels are often very close, especially when we take into account Stark splitting, which could spread each level by several hundred cm  $^{-1}$ . However, a 1.88 $\mu$  absorption line, described as characteristic of U  $^{3+}$  28,29 does not fit into the 5f  $^3$  energy level scheme. Either hydrogenic wave functions are decidedly in error or this 1.88 $\mu$  line belongs to U  $^{4+}$ , where it could fit into the  $^3$ H $_5$  state.

It is difficult to see how these levels can be definitely assigned until the free ion spectra have been analyzed. Also it is quite desirable to know the position of the 6d configuration in order to estimate line strengths. Therefore, a program has been initiated to analyze the free ion spectra of ionized uranium. We hope to start with UIV (U<sup>3+</sup>) and to go on to UIII and UV. This work will be done in cooperation with the Physics Department of The Johns Hopkins University.

CATAR

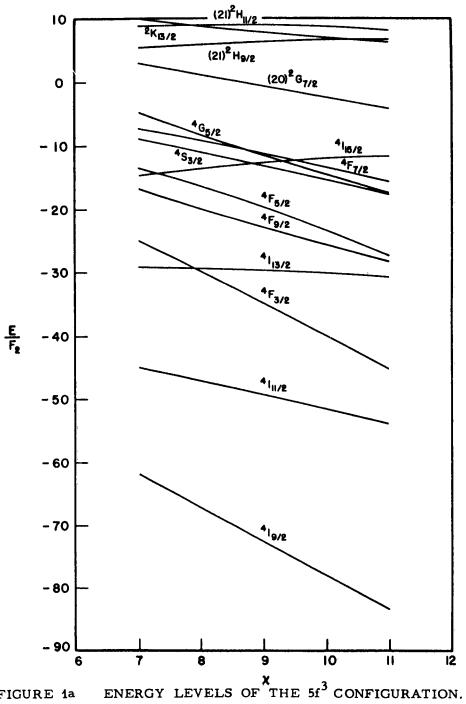


FIGURE 1a

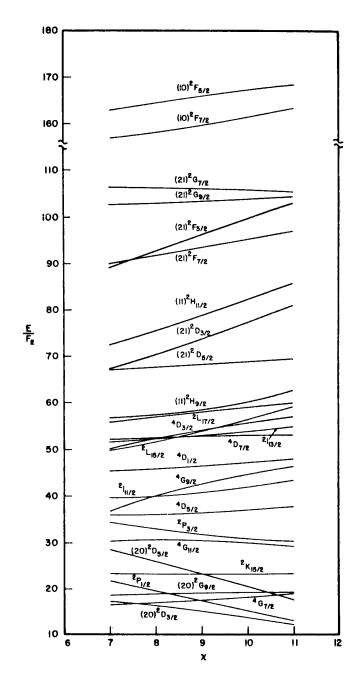


FIGURE 1b ENERGY LEVELS OF THE 5f<sup>3</sup> CONFIGURATION.

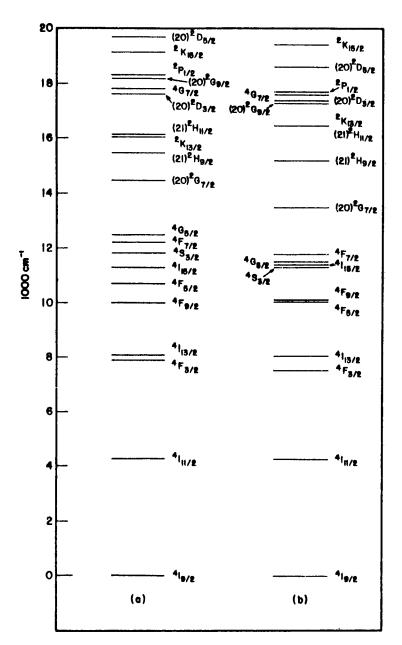
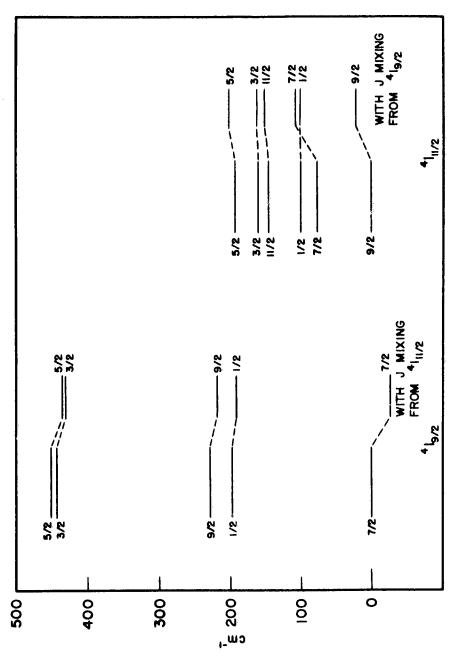


FIGURE 2 LOW-LYING ENERGY LEVELS OF THE 5f<sup>3</sup> CONFIGURATION FOR  $\chi = 8$ .

(a)  $F_2 = 211$   $F_4 = 30.004$   $F_6 = 3.3971$  (hydrogenic)

(b)  $F_2 = 213$   $F_4 = 31.098$   $from U^{4+}$ ).

1 1



CRYSTAL FIELD SPLITTING IN A  $D_{3h}$  SYMMETRY SITE  $A_2^0\langle r^2\rangle=170$   $A_6^0\langle r^6\rangle=-100$   $A_4^0\langle r^4\rangle=-80$   $A_6^0\langle r^6\rangle=1100$ . FIGURE 3

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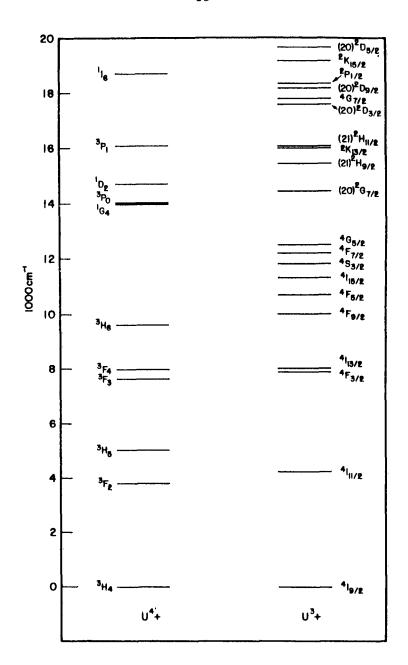


FIGURE 4 LOWER ENERGY LEVELS FOR  $u^{3+}$  AND  $u^{4+}$ .

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EIGENVECTORS

2P 0.76853 -C.63881
4G C.63882 C.76853

REDUCED MATRIX ELEMENTS IN INTERMEDIATE CCUPLING
F2V2 C.CCC C.OCCC
F4V4 C.CCC C.OCCC
F4V4 C.CCCC C.OCCC
G.OCCC
G.OCCC
G.OCCC
G.OCCC
G.OCCC
G.OCCC
G.OCCC
G.OCCC
G.OCCC
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                   EIGENVECTORS

2 0.7254 - 0.68354

4 0.68355 0.72554

4 0.6835 0.72554

FRUCED PATRIX ELEMENTS IN INTERMEDIATE COUPLING

FAV. 0.0000 0.0000

FAV. 0.0000 0.00000

FAV. 0.0000 0.0000

FAV. 0.0000 0.0000
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EIGENVALUES
                                                     EIGENVECTORS
21.714 45.335
EIGENVECTORS
2P 0.22131 -C.57C45
6 0.57C5 C.22131
RECUCEC MATRIX ELEMENTS IN INTERMEDIATE CCUPLING F2V2 C.CCC C.CCC F4V4 C.CCC C.CCC F6V6 C.CCC C.CCC F6V6 C.CCC C.CCC F6V6 C.CCC C.CCC F6V6 C.CCC C.CCC C.CCC F6V6 C.CCC C.CCC C.CCC
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4C 0.62542 0.78636

REDUCEC ##IRIX ELEMENTS IN INTERMEDIATE COUPLING
F2V2 0.6000 0.0000
F4V4 0.0000 0.0000
F4V4 0.0000 0.0000
6 0.40591 0.26674
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                      2C.713 45.586
EIGENVĘCICRS
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                       IS.683 45.866
EIGENVECTORS
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                18.629 46.17C
EIGENVECTCAS
X= 7.C, J= 1/2
EIGENVALUES
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EIGENVALUES
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EIGENVALUES
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EIGENVALUES
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 $5f^{\frac{3}{2}}$ TABLE I. WAVE FUNCTIONS, ENERGIES, AND REDUCED MATRIX ELEMENTS OF

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																										_	_		_	- ·-	
71.821	-0.12654	0.45966	-0.45499		281.5520	00000	1.00165		73-507	,	-0-13754	7380	0.58283		-0.07998	207 9477	0000-0	00000	1.01049		15.262	ATTAL 0- 81326 2	0-47793	0.49537			2 2 2 2 2 2	329-9735	0000*0	1.01855	
53.253	-0.20515	0.47363	0.38575	1.21463 7.1NG	0.0000	0.0000	1.03851		54,125		-0.20483	6.302.0	-0.53413	0.41041	-0.21591	71.1%	070000	000000	1.03395		54.943		10.64-0-	-0.43247	0.55212		G.21/21	252.5767	000000	0,0000	
32.264		0.39419	0.66404	TATE CCUR	3.6662 -38.4231-205.5405.281.5520 0.0000 0.0000 0.0000 0.0000	0.0000	1.09077		31.700		0.24894	0.33194	0-46295	0.63613	0.13475	1141E CEU	-67+8-67-	5000	1.09729		31.227	000	41878 0-	0-43560 -0-43247	0.44857	0.60737	-0.15362	-20-7323-	0.0000	0.0000	2150144
15.855	0.48111		G.43329	INTERFE	3.6662	0000-0	1.16463		15.277				0.35758	-0.46251	0.38824	A INTERPECIALE CEUPLING	. 6000		1.16496		14.643		23C27 3 71EZE OF 71EZE OF 756C7 3	0.37289	-0.38201	-0.45029	-0.38807	-19-7803	0000-0	0.0000	110011
-12.160	C.79618	C.C7631	C.C868C -0.43329	C.36544 -U.38664 -U.II.620 -U.ZI.323 EPENTS IN INTERPEDIATE CCUPLING	99.1254	2220.2	1.62442		-13 227		C. 78383		100000	0.09169 -0.46251	C.38543 -0.38824 -0.13475 -C.21591				1.60624		-14.427		1571.7	70075	. 55500*0-	C.C9629 -0.49029	0.78154	92,4406 110,6751 -19,7803 -20,7323-252,5767 329,9735	2,200-2	0.0000	9
3/2		2P -C.26C28 C.465CC 2C 0.46529 -C.C7651		O.ECGGI FATRIX EL			3 0.61336	C/ E +1	797			15317-0-	(21)20 - 0-4/494 - 0-0120 - 0-40212	4C -C.C9569	0.19344		22 95.9163 105.1022		0	X= 9.5, J= 3/2	-27.359	EIGENVECTORS	784/1*D- 54	7 - C. 28166 C. 1.1064	(21)2C -C.18894 -C.CC499 -0.38201	46 -6-10191	4F C.78154 C.39993 +0.38807 +0.15362 G.21/	2 92-4406		000010 9	2000000
K 3	EIGEN	. 42 (20120	15(12)	4F REDUCED	F2V2	F6V6			EIGEN	EIGEN	4	7	(21)26	. <del>.</del>	4	RECUCED	F2V2	4744		6 HX		EIGEN	•	. 42	(21)2	*	4 (	F2V2	4744	F6V6	
=																															
67.27	0.08990	0.35271	0.69162	-0.03953	176.2528	0000	0.97038		,		-0.10254	0.38344	0.41318	-0.47837	-0.05015		215.5876	0000	0.98175		70-213	:	-0-11482	0.41144	0.63537	-0.46746	-0.06057	250-2035	000000	000000	C0266-0
50.233 67.27	,		,	-0.20527 -0.03953	17 868		0			24.014 00.033	,			٠,	-0.20906 -0.05015		21		1.04869 0.98175		52,318 70,213		ľ			•	00	5			
	-0.20225	C.59253 C.53972	,	121	969		1,05403 0		;		-0.20407	6.57002			-0.08203 -0.20906 -0.05015				0				-0.20497 -	0.54724		•	00	5		0.0000	
50.233	0-13778 -C.20225	-0.27697 C.59253 0.32595 C.53972	,	121	969		1.08088 1.05403 0			51.314	-0.20407	6.57002	0.34869 0.51780	6-71603 6-32753	-0.37734 -0.08203 -0.20906 -0.05015		25.2438 -52.5226-142.0223 215.5876		1.04869 0		52.318		0.19116 -0.26497 -	-0.32109 d.54724	0.3/105 0.49229	0.69074 C.35815 -	00	5		00000 000000	04540
34.471 50.233	0.45368 0.13778 -C.20225	-0.45858 -0.27697 0.59253 0.51321 0.32595 0.53972	-0.2572G 0.49605 -0.43382 -0.34335 0.73992 G.2939G -	121	969		1,15173 1,08088 1,05403 0			13.625 51.514	-0.20407	-0.42854 -0.29986 0.57002	0.34869 0.51780	6-71603 6-32753	906		25.2438 -52.5226-142.0223	0.000 0.0000 0.0000	1.08250 1.04869 0		-11_047 16.383 32.917 52.318		0.19116 -0.26497 -	-0.39772 -0.32109 0.54724	0.3/105 0.49229	0.69074 C.35815 -	00	5	00000 000000 00000	0,000 0,000 0,000	1-68586 1-04546
17-332 34-471 50-233	C.e3843 0.45368 0.13778 -C.20225	C.43417 -0.45858 -0.27697 C.59253 -C.C8175 0.51321 0.32595 C.53972	.0.14578	C.31132 -0.36939 -0.06724 -0.20527	969		1.70728 1.15173 1.08088 1.05403 0			*16.16 66.66 178.61	354 C.82349 0.46667 0.16369 -C.20407	-0.42854 -0.29986 0.57002	0.34869 0.51/80	6-71603 6-32753	4F 0.23312 C.3325C -0.37736 -0.08203 -0.20906 -0.05015	SLEPENTS IN INTERMEDIATE CCUPLING	85.9424 25.2438 -52.5226-142.0223	0,000 0,000 0,000 0,000 0,000	1,15acc 1,0825c 1,04869 0	X# 6.6, J# 3/2	16.383 32.917 52.318		C.aC54C 0.47576 0.19116 -0.2C497 -	-c.2485C C.4563C -0.39772 -0.32109 G.54724	0.3/105 0.49339	0.69074 C.35815 -	00	4 [NIERFELIAIE CLUFLING	00000 00000 00000 00000	000000 000000 000000 000000	I.16233 1.08580 1.04346

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CHERNALISE   STATE   CHERNALISE   CHERNALISE	176.917	1	-0.00952 -0.00952 -0.02593	210.4893 5.000 0.0000	178.325	r	0.71828	-0.01084	-123.5595 202.4244	•								
2 C-1.655	ט ט ט	-0-12401 0-75264 0-31351 0-41229	0.11525	133-5466	163.152	0.13230	0.41218	0.12102 -0.13678	129.8718	1.11136								
2 C-1.655	68-912	C.17764 C.288G6 C.55678	C.24477 - C.24477 - PLING -	-73.0532- 0.0000 1.03911	69.553		C.44115 C.569C7	0.04483 -0.25993 -	3.1252-	1.03276								
2 C-1.655	36.931	0.84941 -0.053&C -0.41174 0.04905	-0.30947 -0.30947 -0.05067 DIATE CEL	17.3089 0.0000 1.20837	37.656	0.85653	0.04162	-0.32838 -0.04604 BIATE CEL	72.6376 29.703d	1.20357								
2 C-1.655	20,575	0.32426 0.57363 0.63984	-0.17216 -0.24681 -4 INTERWEI	149.4069	17-670	0.27970	0.18867	-0.16896 - C.27466 -	-53.0462 156.2565	1.19632								
2 C-1.655		C.18957 -C.C94C1 -C.C7313	C.56513 C.56513 C.767CC LEFENTS 11	351.4905-		C.23354 -C.C9141	C.13345	C.6461C - C.6943Z LEMENTS 13	79,7751 340,6062-	C.81536								
2 C-1.655	,	CTCRS 0.29806 0.02112 0.05229	-0.51896 -0.51896 -26.9667	-50.1219 0.0000 0.50695	LEES 5/2 -27.3CS	C.28461 C.16463	-0.25384	-0.59443 -0.59443	-32,3355	C.E6372								
2 C-1.655	X= 10.0 EIGENV	E1GENVE (20)2D (21)2D (21)2D 4C (10)2F	AFEUCE	7474 7644 6	x= 11.0 EIGENVA	E1GENVE (20)20 (21)20	(10)2F (21)2F	7.4 9.4 0.5 0.5 0.5 0.5 0.5 0.5 0.5 0.5 0.5 0.5	F2V2 F4V4	) (9 )								
2 C-1.655	174.906	0.03420 -0.04239 0.73931	-0.00554 -0.00554 -0.02494	221.2466 0.0000 0.85960	175.563	0.03635	0.73561	-0.02536	188,2085 217,8319 0,0000	0.86000	****	0.03854	-0.05176	0.73136	-0.00746	188-3359	214.2479	0.86045
/ N N N N N N N N N N N N N N N N N N N	94.308	0.10890 0.73957 0.28937 0.41877	0.10544	135.8211 0.0000 1.68859		-0.11434		0.10388	115.2838-	1.09458			0.74923	0.41351	0.11215	116,0287-	134.7696	1.09975
/ N N N N N N N N N N N N N N N N N N N	47,999	72025	0.56062 C.03192 -C.21811 - PLING	79.4345- 0.0000 1.05148	68.296	0.17297	0.43360	0.03481 -0.22759 ·	-77.1691-	1.04688	0	C.17538 .	-C.30414 C.55623	0.43577 C.56580	0.03754 -C.23646	PLING 1.5784-	-75.0485-	1.04279
/ N N N N N N N N N N N N N N N N N N N	36.088	0.81447 -0.12717 -0.48141 0.06803	0.08932 -0.26873 -0.06240 - CLATE CCUR	16.7031 0.000C 1.22058			0.06025	-0.2845C -0.C5764 - CIATE CCUF	63.0676	1.21551	, , ,		-0.07345 - -0.42952	0.05404	-0.29787 -0.05379 -	CIATE CCUF 66.5858	9-7494	
/ N N N N N N N N N N N N N N N N N N N	24.488	0.43345 0.55496 0.66800	-0.18562 - 0.19837 - N INTERME	124.6787 6.0000 1.21192		0.38912	-0.18446 -0.17149	C.21557	135.8950	1.21014		0.35342	0.54966	-0.18698	-0.17515 · C.23166 ·	4 INTERPE	143.7998	1.20747
/ N N N N N N N N N N N N N N N N N N N	ער מ מ נ	C.11489 C.09013 C.0575	C.15967 . C.40291 . C.87C23 . LEPENTS IN	343.3266 0.0000 0.67814	7	13569	C. 16677 C. 15444	C.46CC3 - C.83855 LEMENTS II	45.0600 -	2445		6.16497	10460-0-	C.16C21 -	C.51493 .	LEMENTS 17	352.6663-	C.73291
## 8 8 8 8 8 8 8 8 8 8 8 8 8 8 8 8 8 8	. J= 5/2 Lues	CTCRS C.10522 C.C4466 C.C2728	-C.15841 0.64784 -C.36521 *ATQIX E1	-35.3699	'n	CTCRS C.20576 C.C5645	-0.13219 -0.17582	C.E1371 -G.41937 YATRIX EI	-20.8369 -43.9415 0.0000	0-55418	٠,	CTCAS 0.30299	C.C6876 -	-0.19253	0.77551	#3TRIX E1 -25.4424	-46.8456	C.53C81
	X= 8.5 EIGENVA	E1GENVE (20)20 (21)20 (10)26	RECUCEC F2V2	7474 7474 7474 7474 7474 7474 7474 747	X= 9.C EIGENVA	E1GENVE (2012C (21)2C	(10)2F (21)2F	4.5 4.6 0.8 0.8 0.8 0.8 0.8 0.8 0.8 0.8 0.8 0.8	F2V2 F4V4		X± 9.5 EIGENVA	EIGENVE (2C)2C	121) 150 160	(10)2F (21)2F	4. 4. IT 12	RECUCED F2V2	4744	2 9

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168.8C4 0.13404 0.13404 0.13404 0.01853 0.0185	169.551 0.14170 0.62534 0.62537 0.62537 0.22758 0.22758 0.22758 0.22758 105.2064 44.3821 101.2832	0.14922 0.14922 0.14922 0.13948 0.2326 0.17998 0.23835 0.2808 43.3885 101.6226 101.6226
2 1 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2	N 11 N 4 4 1 H W W W W W W W W W W W W W W W W W W	2 - 10 - 10 - 10 - 10 - 10 - 10 - 10 - 1
106.292 -0.05126 -0.14463 0.25775 0.25775 0.59705 -0.72536 1053.0074	106.195 -0.65861 -0.15431 0.73694 0.59204 0.752101 -107.5418	106.074 -0.06655 -0.16388 0.3232 0.33242 0.58703 0.71660 -0.71660 -109.72865
52.78£ 92.599 106.292 168.8C4 C.96591 C.2C968 -0.05126 0.13404 0.2193 G.2C958 -0.12471 0.155371 C.7622 C.2C059 0.2775 -0.42108 C.7742 -0.2768 0.2775 0.15537 C.7742 -0.2768 0.2775 0.1553 C.7742 -0.2768 0.2775 0.1553 C.7742 -0.2768 0.2775 0.1553 C.7742 -0.2769 0.77536 0.21537 C.7742 -0.2769 0.77536 0.21537 C.7742 -0.2769 0.77536 0.21537 C.7742 -0.2769 0.2775 0.2779 C.7742 -0.2769 0.2779 C.7742 -0.2769 0.2779 C.7742 -0.2769 0.2779 C.7742 -0.2769 0.2779 C.7742 0.2779 C	52.692 93.469 106.195 169.551 0.96317 C.21434 -0.62661 0.1477 0.02965 6.41784 -0.1531 0.6549 0.03965 C.03365 0.31035 -0.40801 0.0354 -0.20942 0.58294 0.16603 0.01554 -0.20942 0.58294 0.16603 0.01554 -C.02988 0.72101 0.22758 0.07996 -C.14476 -0.6861 -0.65416 0.1054 -0.14476 -0.6861 -0.65416 93.0011 -22.027-107.5418-106.2064 13.4776 -42.7447-156.2655 101.2632 13.4776 -42.7447-156.2695 101.2632	52.984 94.332 106.074 170.360 0.96040 0.21834 -0.06655 0.14922 -0.2771 0.4772 -0.1888 0.85514 0.03721 0.1772 -0.1888 0.85514 0.08847 -0.2177 0.1688 0.2324 0.08847 -0.2177 0.58703 0.17298 0.01801 -0.1834 -0.71660 0.2385 0.01860 -0.1834 -0.07763 -0.05808 13.74F CUPLING 73.0536 -32.0117-109.2285-102.5593 27.2811 -42.5152 0.41386 13.7645 -0.4514-158.5973 10.6225 1.40331 1.13921 0.42552 1.12670
17.439 52.78¢ 92.599 106.292 168.80¢  C.C321¢ C.96591 C.20968 -0.05126 0.1340¢  0.07154 -0.07193 C.4839 -0.14463 C.85432  0.07554 -0.0745 -0.02371 C.00637 0.04771 C.15677  0.4657 0.1325 -0.02059 0.7255 -0.21537  0.4677 0.1325 -0.02059 0.72538 0.21637  0.4677 0.1325 -0.02059 0.72538 0.21637  10.4677 0.1325 -0.02059 0.72538 0.21637  10.4677 0.1325 -0.02059 0.72538 0.21637  10.4687 0.1325 -0.02059 0.72538 0.21637  10.4687 0.205 -0.02059 0.72538 0.21637  10.4687 0.205 -0.02059 0.72538 0.21637  10.4687 0.205 -0.02059 0.72538 0.21637  10.4687 0.205 -0.02059 0.72538 0.21637  10.4687 0.205 -0.02059 0.205 0.205 0.205  10.4687 0.205 0.205 0.205 0.205  10.4687 0.205 0.205 0.205 0.205  10.4687 0.205 0.205 0.205 0.205  10.4687 0.205 0.205 0.205 0.205  10.4687 0.205 0.205 0.205  10.4687 0.205 0.205  10.4687 0.205 0.205  10.4687 0.205 0.205  10.4687 0.205 0.205  10.4687 0.205 0.205  10.4687 0.205  10.4688 0.205  10.205  10.205  10.	17.766 52.692 93.469 106.195 169.551 0.03405 0.98317 0.28434 -0.05861 0.14170 0.07142 -0.05645 0.83345 0.31035 -0.40801 0.57546 -0.05655 0.83345 0.31035 -0.40801 0.55354 0.020942 0.52094 0.16603 0.165036 0.056354 0.02037 0.46672 0.01554 -0.09883 0.72101 0.22755 0.14059 0.166031 0.16603 0.11603	18.1C8 52.984 94.332 106.C74 17C.360 -0.03595 0.9604C C.21834 -0.C6655 G.14922 0.0724 -0.09521 G.89410 0.32242 -0.39421 0.52325 -0.03721 C.16266 0.73912 0.62226 0.56070 0.08847 -C.21771 0.58703 0.17298 0.56070 0.08847 -C.21771 0.58703 0.17298 0.4528 0.01801 -0.1834 -0.07166 0.23825 0.45816 0.02846 -0.1834 -0.07763 -0.05808 55.633 -93.0536 -22.117-109.7285-102.5593 35.8252 27.2811 -42.5152 27.1260 0.3382 1.00538 1.40331 1.13921 0.42552 1.12670
17.439 -6.0214 0.07163 - 0.07549 - 0.6695 0.6695 0.65513 N. IMTERPE 36.2557 14.6892	17.766 52.892 0.03405 0.96317 0.07492 0.01956 0.077492 0.01956 0.56756 0.02854 0.56376 0.02854 0.56376 0.02854 0.56376 0.01854 0.66072 0.01854 0.1778PEDIATE CC 55.46672 0.01956 0.1778PEDIATE CC 55.46672 0.01956 0.107988 12.7557 10.3888 12.4776	18.1C8 -0.03595 0.07229 0.07229 0.05225 0.05225 0.052225 0.052225 0.052225 0.052225 0.052225 0.052225 0.052225 0.052225 0.052225 0.052225 0.0525 0.0525 0.0525 0.0525 0.0525 0.0525 0.0525 0.0525 0.0525 0.0525 0.0525 0.0525 0
' ' 🛱	' ' A	ELGENALUES  ELGENALUES  ELGENVELTES  LIC.434 — L.42E 18.1CB 52.984 94  4. G.CL639 — C.C414E — 0.03595 0.9604C C.211  (10.72F — C.CE527 C.13225 C.07229 — 0.22771 C.407  4. C.CC6059 C.15042 0.07229 — 0.23771 C.407  4. C.CC6059 C.15042 0.07229 — 0.03721 C.407  (20.12G — 0.49598 — C.E99315 0.55070 0.08847 — C.C217  4. G.C.2324 C.148.2 — 0.45280 0.02847 C.2217  (21.12G — 0.49598 — C.E9376 0.45280 0.02847 — C.C217  (22.12G — 4.48.12 — 0.45280 0.02847 — C.117  FEVE — F. S.
ELGENALUES  ELGENALUES  ELGANALUES  ELGANALUES  ELGANALUES  ELGANALUES  ELGANALUES  ELGANALUES  (21)2F -C.C377  (21)2F -C.C377  (21)2F -C.C377  (21)2F -C.C377  (21)2F -C.C377  (21)2F -C.C377  (21)2G -C.439  (21)2F -C.438  (21)2F -C	EGGENVALUES EIGENVALUES EIGENVALUES EIGENVECTERS -C.522 EIGENVECTERS -C.6311 EIGNES -C.6312 EIGNES -C.6313 EIGNES -C.6313 EIGNES -C.6313 EIGNES -C.6313 EIGNES -C.6313 EIGNES -C.6313 EIGNES -C.6314 EIGN	ELGENVALUES ELGENVALUES ELGENVECTERS 4C G.CL629 -C.C414E (10.12F -G.C827 C.15425 (12.12F -G.C827 C.15425 (12.12F -G.C827 C.15425 (12.12F -G.C827 C.16425 (12.12F -G.C827 C.12F -G.12F
EIGENATURES EIGENATURES EIGENATURES (10) 2F -0.05 (10) 2F -0.05 (20) 2G -0.40 (20) 2G	EIGENVLUES EIGENVLUES EIGENVECTURS EIGENVECTURS (10.5	EIGENATURES EIGENA
w 2-1001004 7-1007	ថ្ងៃសំសុស្សស្ត្រីក្រុង	p 487.50 NW N4WF
166.943 0.11037 0.85141 0.45898 0.13423 0.1804C 0.C3894 16.2887 49.2103 96.4088	167.499 0.11836 0.85262 0.61665 0.19275 0.19275 0.19275 0.19275 0.19275 0.19275	168-119 0-12626 0-85358 0-643397 0-01676 0-15690 0-20475 0-204643 1-13067
106.426 166.943 0.02254 0.11037 0.11516 0.85141 0.2278 -0.01341 0.6153 0.1342 0.73756 0.11804 0.6523 -0.1804 0.6523 -0.1804 0.6523 -0.2084 97.9995-110.2687 43.8055 4.2103 43.7133 94.4088	106.409 167.499 0.03924 0.11836 0.12633 0.85282 0.03016 0.01595 0.03016 0.01595 0.03874 0.01477 0.73367 0.11677 0.04499-113.4673 46.3260 0.4499-113.4673 47.3996 98.1541 0.91258 1.13198	106.364 168.119 -0.05448 0.12626 -0.1367 0.83538 -0.63245 0.01676 -0.63245 0.01676 -0.66215 -0.0463 -0.06215 -0.0463 -0.06215 -0.0463 -0.06215 -0.0463 -0.06215 -0.0463
66 106.426 43 -0.03254 9940.11516 60 0.25278 60 0.25278 60 0.77756 776 0.05523 776 -0.05523 777 -0.05523 778 -0.05523 778 -0.05523 778 -0.05523 778 -0.05523 778 -0.05523	146 106.409 226 -0.03924 83 0.2833 83 0.2833 83 0.2016 51 0.2834 117 -0.03874 117 -0.03874 121 43.256 10-147.3996	91.724 106.364 168.119 C.2C433 -0.C448 0.12626 G.43382 -0.1347 0.85358 C.4316 0.03245 0.01676 C.1917 0.03245 0.01676 C.1917 0.03245 0.01676 C.1917 0.06215 -0.0443 J.1146 30.9428-102.8838-110.6855 38.782-102.6849 9.543
66 106.426 43 -0.03254 9940.11516 60 0.25278 60 0.25278 60 0.77756 776 0.05523 776 -0.05523 777 -0.05523 778 -0.05523 778 -0.05523 778 -0.05523 778 -0.05523 778 -0.05523	146 106.409 226 -0.03924 83 0.2833 83 0.2833 83 0.2016 51 0.2834 117 -0.03874 117 -0.03874 121 43.256 10-147.3996	52.666 91.724 106.364 168.119 0.56863 C.2C433 -0.05448 0.12626 -0.0306 G.4382 -0.1347 0.65358 -0.0396 C.8306 0.4387 0.65358 0.0173 0.09259 0.03245 0.01676 0.00115 -0.0877 0.7296 0.20475 0.00185 -0.1841 -0.06215 -0.06443 124.74 C.UPLING 12.521 46.8512-106.6819 46.554 11.5792 46.8512-106.6840 9.5433 11.40905 1.14139 0.91568 1.13067
16.571 52.382 89.966 106.426 0.02632 0.9740 0.459940.11516 0.07103 -0.02735 0.03838 0.22578 -0.24774 0.02736 0.04573 -0.02735 0.03838 0.02578 -0.15716 0.04633 0.00752 -0.07776 0.02736 0.5772 0.05637 0.07776 0.02736 1.5772 0.05637 0.05777 0.02737 53.496 -92.409 -22.999 -97.99957 23.425 29.9339 -50.3337-143.7733 0.528392 1.4124 1.14254 0.9992	6.839 52.532 90.846 106.409 02829 0.97132 C.19926 -0.03924 07073 -0.19493 C.4945 -0.12303 070764 -0.02937 C.49328 0.03016 4507 -0.02960 C.6993 0.03016 4507 -0.02960 C.6993 0.03016 4504 0.00924 -0.03151 0.05834 4804 0.00924 -0.03151 0.05834 4804 0.00924 -0.03151 0.05834 4805 -92.874 -93.4254 4603 229.3955 -37.3421 48.2260 45603 1.41087 1.14200 0.91258	24 106.364 33 -0.6448 92 -0.13487 66 0.03245 995 0.03245 774 0.7266 71 -0.66215 726-102.8938-10388 122-150.6640
16.571 52.382 89.966 106.426 0.02632 0.9740 0.459940.11516 0.07103 -0.02735 0.03838 0.22578 -0.24774 0.02736 0.04573 -0.02735 0.03838 0.02578 -0.15716 0.04633 0.00752 -0.07776 0.02736 0.5772 0.05637 0.07776 0.02736 1.5772 0.05637 0.05777 0.02737 53.496 -92.409 -22.999 -97.99957 23.425 29.9339 -50.3337-143.7733 0.528392 1.4124 1.14254 0.9992	6.839 52.532 90.846 106.409 02829 0.97132 C.19926 -0.03924 07073 -0.19493 C.4945 -0.12303 070764 -0.02937 C.49328 0.03016 4507 -0.02960 C.6993 0.03016 4507 -0.02960 C.6993 0.03016 4504 0.00924 -0.03151 0.05834 4804 0.00924 -0.03151 0.05834 4804 0.00924 -0.03151 0.05834 4805 -92.874 -93.4254 4603 229.3955 -37.3421 48.2260 45603 1.41087 1.14200 0.91258	1.254 17.129 52.666 91.724 106.364 168.119 -C.03125 -0.03023 0.56863 C.22433 -0.54448 0.12626 -C.103129 0.07129 -0.03030 C.43892 -0.13467 0.58358 -C.10317 0.07581 -0.03936 C.43937 0.03245 0.03346 0.643937 -C.46625 C.47157 -0.03173 0.09295 0.03324 0.01676 -C.163715 0.05893 0.03115 0.09295 0.03245 0.01676 -C.184625 0.47157 0.01115 -0.08774 0.72940 0.020478 -C.184625 0.47157 0.01115 -0.08774 0.72940 0.020478 -E.84715 11 11 11 11 11 11 11 11 11 11 11 11 1
2 2.582 16.571 52.382 89.966 106.426 -C.C244E -0.02632 0.9740C C.19143 -0.02254 C.C2456 0.077008 -0.18587 0.459940.11516 -C.4564 0.4776 0.02735 0.05954 0.0.2586 -C.4564 0.4776 0.02735 0.05954 0.0.25878 -C.2464 -0.44776 0.02735 0.05954 0.0.25878 -C.2464 -0.446933 0.05757 0.05954 0.0.2986 -C.2164 -0.46933 0.05757 0.059591 -0.17274 -0.05523 -0.2986 -2.2466 -0.46933 0.05959 -0.2996 97.96995-1-294,226 33.4699 -0.29895-1-284,226 33.46995-1-284,226 34.2699 -0.29895-1-284,226 34.28995-1-284,226 34.28995-1-284,226 34.28995-1-28895-1-28	6.839 52.532 90.846 106.409 02829 0.97132 C.19926 -0.03924 07073 -0.19493 C.4945 -0.12303 070764 -0.02937 C.49328 0.03016 4507 -0.02960 C.6993 0.03016 4507 -0.02960 C.6993 0.03016 4504 0.00924 -0.03151 0.05834 4804 0.00924 -0.03151 0.05834 4804 0.00924 -0.03151 0.05834 4805 -92.874 -93.4254 4603 229.3955 -37.3421 48.2260 45603 1.41087 1.14200 0.91258	.ues .ues
2.562 16.571 52.382 89.966 106.426 C.C244E -0.C2632 0.9740	16.839 52.532 90.846 106.409 -0.02829 0.97132 0.19926 -0.03924 0.07073 -0.19939 0.4967 -0.12503 0.07614 -0.02937 0.49283 0.26837 0.45694 0.00924 0.69833 0.03016 0.45694 0.00924 -0.03151 0.65685 -0.4594 0.00924 -0.03151 0.05874 -0.48940 0.06686 -0.4694 0.20924 -0.3151 0.05874 -0.48940 0.06686 -0.4603 29.3395 -37.3421 43.2260 24.346 12.0029 -48.6710-17.3996	2 1.254 -C.C3125 -C.12373 C.12373 C.12373 C.12375 C.16661 -C.16661

TABLE I CONTINUED.

X= 8.5, J= 9/2						X* 10.00 J* 9/2	
E14ENVALUES -21.232	12 6-241	19.010	41.451	57.779	103.181	935 -25.516 6.768 19.257 44.779 50.074	103.874
CTCRS					1	CTERS	
C.C2624 -C.		-0-C1652	0.48698	0.13730	0460-0	10-53616 0-11461 0-04148 0-41340 0-23249	47.47.0
-6.C8043 C.4	1	66240-0	C749C-0	6,971-0	0.04040	445270 077150 FIRET C-37150 0 78780 0	0.6400
		2,067.0			1521	0 053001 -0 36145 -0 32530 0 88588 -0 05460 0 30780 -	17170
	14/91-D-		20000		# # # # # # # # # # # # # # # # # # #	C. 10414 -	100000
	20020-0-	-0-887-0-	CERS7-0-		98447-0	00477-0 34134-0-36361-0-36361-0-36340-0-36361-0-36661-0-36661-0-36661-0-36661-0-36661-0-36661-0-36661-0-36661-0-36601-0-36601-0-36661-0-36601-0-36601-0-36601-0-36601-0-36601-0-36601-0-36601-	67007.0
(21)2F -C.367E3 C.57C1E	16764-D 31	16662-0		1000	1120-0	CODE OF FIRE OF CALL OF CARE OF ACTION OF	10000
41 0.51406 0.349	14 0-16833	0.09232	£		C+700-0		76001
RECUCED WATELY ELEMENTS IN INTERPECTATE COUPLING	IN INTERPE	מזאוב כרס	1				
F2v2 -28.27C4 37.36	18.4934	0.2187	701		B990.09-	3-1-0236 IU-2200 3-3312 -II-9653 -I-6163 -6	-61.6336
2	3C -18.9634	11-1594	13.6331	-1-7356	3.5128	12.10/5 13.6854 -1.5442	1654.2
		24.9460	2.4681		-78.8337	85.6240 -45.4928 67.7099 36.4706 -19.0664	-81.4340
G C.76C42 1.C17	rc 1.18971	1.12938	1.10453	10,7407	1.09790	1.11682 1.07219 1.02020	1.05393
x= 9.C, 0= 9/2						X= 11.C+ 9/2	
i			***	746 03	102 207	377 67 676 77 18 718 77 301	357 701
1/0-77- 242-71-71-71-71-71-71-71-71-71-71-71-71-71-	0 1	14-67	171.75	10.00	100*001	7 TT - 17 TT -	
4F C.C2922 -C.33622	22 0.78621	0.00356	G-48562	0.16755	0.06434	549 -C.1362C G.7606C G.C9597 C.45297 G.29229	0.08594
	•	0.04994	C.54935	0.15868	0.64627	-C.2671C C.C674C C.46663 0.27715	C.45209
٠,				-0-45560	0.70605	-C.38365 C.07269 O.11224 -C.19559 -0.57459	6.67635
C.C4726 -C.	٠,				-0.05039	0.C5818 -C.2675E -C.26832 0.8622C -C.12148 0.3C37C	-C.C5134
C-12334	16 -C.07057	-0.30421	-0.35317	0.80394	0.26287	-0.00526 -0.35855 -0.54984 0.64222	0.31298
	34 0.49772	0.25966		0.12318	0.08977	C.52387 C.49056 0.29296 -C.46539 -0.C2695	08963*?
41 0.50325 0.36021		494L3-0- 84863.0		-0.C2294	0.00133	-0.04695	.00023
RECUCED MAIRIX ELEMENT		CIATE CCU					
7	.5577 15.5035	1.9570	261		-63.5283	36.4272 3.5794 10.2791 -28.0794 13.1508	-59.2839
F444 -16,C655 2,6276	76 -19.1481	11.5072			3.2148	2.0133 -19.6496 12.5202 12.4803 -0.4163	1.4665
52	30 -41.6033	59.3157	13.7014		1669*61-	79.2552 -48.4198 75.312C 54.3491 -40.1685	-53,2367
G 0.76275 1.C1748		1.12466	1.09483	0.98819	1.09666	G 0.77149 1.C155E 1.18305 1.1088C 1.C5C8C 1.C5137 1.	1.09123
C/25 HT 10 - C/25							
-75,229 -24,059	55 6.617	19-178	43.829	59.137	103-618		
-0- 610		0.02496	0.48097	0.19984	0.06932		
(20)26 -C.C3931 C.43657	,			0.19159	0.64733		
(21)2G C.C8338 -C.375Gl	0.05010	0.10122	-C-31867 -	-0.49029	0.69936		
ij	-0.26497			0.28925	-0.05103		
(11)2F 0.12597 -C.21317	-0.05610	-0.31356		0.76888	0.27567		
(21)2h -0.38874 C.55C36		0.26666		0.08540	0.09207		
4I C.5C255 C.37C&I	51 0.17480	0-10494 -0-07473		-0.62940	91100-0		
AECUCED MATRIX SLEPENT	ENTS IN INTERPECIATE CCUPLING	CIATE CCU					
-	13.1966	2450-5		. 9501.01-	2001-20-		
7 :		11.6260	13.9360		46 642		
F6V6-132.1015 38.9322	1	2400.00	1 00530		1 09527		
:	1.186/8	1.16017	100001	10000	17660-1		

101.77	C.32119 C.66437	C.47989	C.04863	JPLING -3 3144		60.4979			78.743		0.33551	C.65252	C.48656	0 7 1 20 0	PLIN6	-3.7653	13.5016	59.6235			8C.419		0-34916	0.64099	13654-0	70774	0.04030 0.136	-3.7144	14-1854	58.7132
40.344	0.66067	0-19949	0.10688 -0.0109C	-4-5479		-1.2879 -17.9907 1.09706 1.08280			4C.8CS		0.69336	-0-01057	0.49114 0.40014	0.11451 ±0 CORE1	CIATE CCU	-3.8436	23.9323	1.5033 -18.9089 07944 1.09829			41.366		0.71525	-0.06411	0.20988	2000000	CIATE COL	-3.2857	22.5906	3.6309 -19.227C
30.532	0.47385	0.22724	G.1C688	N INIEKFE -4-4371	11.8203	-1.2879 -			30.347		0.40003	-6-73556 -0-61657	C-23203	0.11451	N INTERPE	-5-C176	11,1525	1.07944			30.088		0.33211	0-16113 -C-13255 -0-74809 -0-06411	0.574034 U.2U988	0-12616 -0-0063	MATRIX OFFERENTS IN INTERSEDIATE COLD INC	-5-4267	10.4588	3.6309
9.150	-C.4844C -C.18675	C.79813 -C.23864	C.19113	C. 1795		-3C.3862 1.11942			5*II		Cathory Itaayaas	70007 * J		0.19305	LEFENTS I	-C.0115	8.2627	-31.538C 1-11937			5.022	,	10.504.0	- 55257-0-	100000	6-19414	LEPENIS II	-C.2294	9.5355	-32.5658
EIGENVALUES -46.101	Elbenvelilks 4G C.C3441 (11)2H O.C8271	(21)2h -C.2C317 C.79E13 . 2I -C.C3363 -C.23864	4I 0.57446 C.19113 G.1C688 -G.C109C C.04	F2V2 -3.8168	-16.4305	F6V6 -15.9618 G 0.57141	9.C. J= 11/2	EIGENVALUES	-45.238	EIGENVECTURS		,	21 -C - C347C - C343C	4I 0.57286	u	F2V2 -3.7584	F4V4 -1C.35C1	F6V6 -15.8U85 -31.538C G 0.57182 1.11937		X= 9.5; J= 11/2	-5C.38C	9	0/404.0- 64580.0 45(ff)	J	71 -0 -1257 -0 -17	4I 0.5713C	٠,	FZV2 -3.78C6	F4V4 -1C.2718	F6V6 -15-66C0 -32-585E
72.464			35	88	9489	1741	# <b>.</b>		73.929					249	!		1.33	) C)		# X						559		618	963	2 6 6
			98 C.03935 CCUPLING	47 -3.9488		99 62.7741 62 1.06748				59 C.29058	-		_	74 C-04249	COUPLING		10 42 Capp				15 75.496	16408-0 81	_			C C-04559	CUPL ING		7 61 3360	
39.321	0.65939 -0.48852 -0.61274 -0.24711	7	0.CZ598 EDIATE CCI	-7-1947	28.2027	-11.6099			39.823	-0.55659		-0-180	0.78945	0.02174	EDIATE	-6.3023	-14.176	1.04258			40-005	-0.61508	-0-1164	-0.1909	0.75589	0.01650	DIATE C	-5.3870	7265 91-	77590
30.326	96659	0.27720	0.07783 I INTERME	-2.0331	14.5828	-13.3387 1.16618			30,553	0.62845	-0.65612	0.22705 -0.18062	0.33964	C.C8747	Ζ.	12.8033	7021-41- 4215-8-	1.14239			30.613	0.55144	-0.69114 -0.11644	0.22589 -0.19093	0.35720	0.09721	MATRIX ELEMENTS IN INTERMEDIATE COUPLING	-3.6986	75-79-01-07-17-17-17-17-17-17-17-17-17-17-17-17-17	1,11853
	- 1		ິຊ											111	٠.	٠,													٠.	ı
.UES -44.734 8.959 30. TCRS	- 1	-C-17713	C.17982 ELEPENTS IN	-3.8749 C.6147	F4V4 -1C.6834 3.8471	c+57613 1-11724 1-16618 1-62162	X= 7.5, J= 11/2		390°5	C.C2998 -C.45817	-C-23432	(21)2h -6.15011 C.81355	-C.19652	0.57778 C.1844E	FRIKIX ELEMENTS	7,44,7	-2C.2821 -27.8527	1.11831	XX 8.0. 13.7.		5.133	0.03222 -0.47200	-0.21135		-C.C3250 -C.217C4	C.18826	LEMENTS I	5.3463 8.010 8.010	-29.1512	1.11965

```
EIGENVECTORS

21 -0.7514 -0.19383 0.97314

21 -0.7515 -0.21446 0.03347

22 0.20345 0.55726 0.20335

RECUCED MAIRIX ELEMENTS IN INTERMEDIATE CCUPLING
FAVA -3.3937 -56.2375 -37.3045
FAV4 -3.3937 -56.0645 1.10706

G 1.10C28 0.94476 1.07090
                                                                                                                                                                                                                                                                                                            52,336
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                   -25.216 8.531
EIGENVECTORS
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                X= 8+C, J= 13/2
EIGENVALUES
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                               X= 8.5, J= 13/2
EIGENVALLES
                                                                                                                                                                                                                                                                               X= 7.5, J= 13/2
EIGENVALUES
X= 7.C, J= 13/2
EIGENVALUES
                                                                                                                                                                                                                                                                                                                                                 EIGENVECTORS

EIGENVECTORS

EIGENVECTORS

EIGENVECTORS

(1124 0.05436 0.76439 0.10461 0.38612

(21)2H -0.2595C 0.74847 0.28359 0.10463 0.60332

(21)2H -0.2595C 0.74847 0.28359 0.10463 0.60362

21 -0.13847 -0.3563 -0.53241 0.60376 0.46460

41 0.5667 0.13663 -0.15565 0.01036 0.60468

RECUCEC PARIX ELEMENTS IN INTERPEDIATE COUPLING

FAVZ -2.73C2 -1.0355 -5.3216 -2.2926 -3.5567

FAVK -10.0493 13.5891 7.9357 18.9502 16.0731

FAVK -10.0493 13.603 7.0807 -17.935 5.8076

G 0.57335 1.11463 1.04516 1.12322 1.08129
                                                                    46 C.C4CE3 -C.51248 +0.27C47 -0.72898 C.36213
(11)2h C.C2CC4 -C.10414 0.77647 0.11042 C.62578
(21)2h -C.23CC4 -C.17CC4 -0.25237 -0.21226 C.46375
21 -C.23667 -C.2CE45 -0.52195 0.64135 C.46875
41 0.56477 C.19425 -0.51395 0.64135 C.46875
RECUCEC PATRIX ELEMENTS IN INTERMEDIATE CLOUPLING
F2V2 -3.7633 -C.4734 -5.6798 -2.8578 -3.6625
F4V4 -1C.1956 10.8554 9.6943 21.298C 14.8427
F6V6 -15.5161 -3.55CE 5.2038 -19.0361 57.7777
                                     41.986
                                   29.819
                                                                                                                                                                                                                                                                                                                    X= 11.C, J= 11/2
EIGENVALUES
 x= 1C.C, J= 11/2
EIGENVALUES
                                 -51.529
EIGENVECTORS
```

EIGENVECTORS

EIGENVECTORS

21 +0.09714 -C.21147 0.97255

41 0.5662 -C.27555 0.03604

2K C.26637 C.29774 0.22591

RECUCED MATRIX ELEMENTS IN INTERMEDIATE CCUPLING
FAVA -2.3924 -4.7632 4.6588
F6V6 -3.3736 1.661 0.5245

G 1.09558 C.55259 1.06937

1

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1

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, L

CATH

```
2K C.48468 -C.5457 0.62285 0.00000

2K C.48468 -C.54657 0.62285 0.00000

15 2L -0.18258 0.70021 0.600000

17 2L 0.000000 0.000000

RECUCED MATRIX ELEMENTS IN INTERMEDIATE CCUPLING

FAV2 -12.0864 -27.0616 -27.4826 -19.6078

F4V4 -2.5312 -C.65538 -4.0736 -3.5651

F6V6 -22.0022 -8.0816 12.5327 1.6876

6 1.16005 1.03323 1.01456 1.05900
                                                                                                                                                                                                                                                                                                                                     6C.053
                                   59.053
                                                                                                                                                                                                                                                                                                                                     59.060
                                     56.358
                                     23.169
                                                                                                                                                                                                                                                                                                                                     23.156
                                                                                                                                                                                                                                                                                                                               -11.699
EIGENVECTERS
                                                                                                                                                                                                                                                                                                  X= 11_C, J= 15/2
EIGENVALUES
                               -11.97C
EIGENVECTORS
X= 10.0; J= 15/2
EIGENVALUES
                                                                  2K 0.41560 C.65635 0.6296 0.0000C
2K 0.41560 C.65635 0.62969 0.0000C
15 2L 0.12622 C.64349 0.7555 0.0000C
17 2L C.CCCC C.CCCC 0.0000C 1.0000C
17 2L C.CCCC -C.CCCC 0.0000C 1.0000C
F2V2 -1C.6269 -26.755 -29.2478 -10.6078
F4V4 -2.6259 -6.795 -3.5651
F6V4 -2.2.8641 -7.7638 14.0767 1.6876
6 1.17285 1.03522 0.99977 1.05900
                                                                                                                                                                                                                                                                                                                                                                   2K 0.43962 C.61554 -0.2033C 0.0000C
2K 0.43962 C.6159C 0.64999 0.0000C
15 2L -0.14466 -0.46659 0.7325 0.0000C
17 2L 0.0000C 0.0000C 0.0000C
18 EUCEC PATRIX ELEPENTS IN INTERPEDIATE CUPLING F2V2 -11.0797 -26.922 -28.6307 -19.6078 F4V4 -2.2625 -0.7259 -3.9321 -3.5651 F4V6 -2.2625 -7.8872 13.6013 1.6876 6 1.16886 1.03405 1.06499 1.05900
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                     2X 0.46270 -C.55324 0.66764 0.00000

2X 0.46270 -C.55324 0.66764 0.000000

15 2L -0.16250 0.66674 0.71074 0.000000

17 2L 0.00000 0.000000 1.00000

FEVZ -11.5670 -Z.7.0216 -Z.8.0421 -10.4678

F4V4 -Z.5659 -G.66210 -4.0097 -3.5651

F4V6 -Z.2.4632 -7.9910 13.6831 1.6876

G 1.16459 1.03343 1.00993 1.05900
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                      58.553
                                                                                                                                                                                                                                                                                                                                     58.053
                                                                                                                                                                                                                                                                                                                                     53.908
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                      55.098
                                     52.791
                                                                                                                                                                                                                                                                                                                  -12.564 23.213
EIGENVECTORS
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                  -12.227 23.185
EIGENVECTCRS
                                     23.246
                           -12.982
EIGENVECTORS
X= 8.5. J= 15/2
EIGENVALUES
                                                                                                                                                                                                                                                                                                  X≠ 9.C, J≠ 15/2
EIGENVALUES
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                  X= 9.5, J= 15/2
EIGENVALUES
```

TABLE I CONTINUED

j

1

.

									M					
J	k	₫₹	1000 f	1/2	3/2	5/2	7/2	9/2	11/2	13/2	15/2	17/2	19/2	21/2
3/2	2	4.5	223.606 798	-1	1									
5/2	24	4.3.5.7 36.7	48.795 005 62.994 079	-4 2	-1 -3	5								
7/2	246	4.2.3.5.7 36.2.7.11 4.2.3.11.13	34.503 279 13.430 383 17.069 718	-5 9 -5	-3 -3 9	-13 -5	7 7 1							
9/2	246	4.3.5.11 36.5.11.13 4.3.5.11.13	38.924 947 6.232 980 10.795 838	-4 18 -8	-3 3 6	-1 -17 10	-22 -11	6 18 3						•
11/2	246	4.3.5.7.11.13 36.2.7.11.13 4.2.3.11.13.17	4.080 444 3.724 918 4.140 015	-35 28 -20	-29 12 4	-17 -13 25	-33 11	25 -27 -31	55 33 11	:		ľ		,
13/2	246	4.2.5.7.13 36.2.7.11.13.17 4.5.7.11.13.17.19	16.574 839 0.903 425 0.393 248	-8 108 -200	-7 63 -25	-13 185	-2 -92 227	-132 -11	-77 -319	13 143 143				
15/2	246	16.3.5.7.17 144.7.13.17.19 16.3.5.13.17.19	5.917 263 0.486 068 0.996 142	-21 189 -75	-19 129 -25	-15 23 45	-9 -101 87	-1 -201 59	-221 -39	21 -91 -117	273 273 65			
17/2	246	36.2.5.17.19 36.2.11.17.19 36.11.13.17.19.23	2.932 564 1.977 134 0.161 702	-40 44 -440	-37 33 -209	-31 13 145	-22 -12 439	-10 -36 481	-51 169	23 -47 -377	-12 -650	68 68 442		
19/2	OF ID	4.3.5.7.11.19 36.2.5.7.11.17.19.23 100.2.3.11.13.17.19.23	3.375 221 0.069 848 0.039 609	-33 1188 -1716	-31 948 -988	-27 503 195	21 77 1353	-13 -687 1931	-1187 1497	-140 <u>2</u> 6	23 -1122 -1870	39 -102 -2346	57 1938 1938	
21/2	246	4.5.7.11.23 36.5.7.11.13.19.23 4.3.5.11.13.17.19.23	5.313 439 0.112 695 0.125 254	-20 702 -520	-19 585 -338	-17 365. -30	-14 70 303	-10 -258 537	-5 -563 558	-775 306	-810 -170	16 -570 -646	25 57 -646	1197 646

(a) S<sub>K</sub>O

J	7/2,-5/2	9/2,-3/2	11/2,-1/2	13/2, 1/2	15/2, 3/2	17/2, 5/2	19/2, 7/2	21/2, 9/2
7/2 9/2 11/2 13/2 15/2 17/2 19/2 21/8	1179.685 678	1.825 742 18.956 090 7.236 272 42.878 566 168.267 644	1.023 533 13.403 980 5.693 469 35.874 782 146.458 185 45.796 652	6.904 105 3.765 875 26.739 484 117.166 548 38.384 024	1.837 559 16.911 535 84.095 184 29.732 138	7.946 248 51.497 573 20.752 510	23.531 134 12.401 997	5.546 342

(b) S<sub>6</sub>

Table II. Stevens coefficients  $\mbox{additional tables of } S^q_K \mbox{ May be found in reference 2. }$ 

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